

> From: Margaret Spence <mspence@parametrix.com>  
> Date: Thursday, January 11, 2007 1:39 pm

> Subject: Re: A few modifications to the table

>

> > Hi gang. I'm working at home again today and will be joining in the

> > conference call. I've attached an Excel spreadsheet (and a PDF

> > print-out of it) I worked up yesterday to guide me through the

> > analysis process and keep files, etc. organized. It also includes

> > several questions I came up with yesterday as I started working

> > through QM.

> > Hopefully, these can get resolved during today's call.

>

> > If anybody needs to reach me, call my cell phone at (b) (6).

>

> Ben, (b) (6) !

>

> Margaret

>

> Margaret Spence

> > Phone: 425-458-6369

> > Fax: 425-458-6363

> > mspence@parametrix.com

>

> PARAMETRIX

> > Inspired people - Inspired solutions - Making a difference

>

> >>> <Benjamin.Shorr@noaa.gov> 01/10/07 10:31 PM >>>

> Sounds good-

>

> > I am trying to use the spreadsheet as a guide- added a couple of

> > fields for Cumulative Distr. charts, summary by areas graphs & maps

> > and am

> > checking them off as I go...

>

> (b) (6)

> Perhaps tomorrow after the call and discussion of progress we

> > can see

> > if I need to spread some pieces that I am responsible for to

> Margaret, > Jim or Carrie to ensure that they get enough attention.

>

> > Thanks,

> > Ben

>

> > ----- Original Message -----

> > From: Robert Gensemer <rgensemer@parametrix.com>

> > Date: Wednesday, January 10, 2007 6:14 pm

> > Subject: Re: A few modifications to the table

>

> > > I think we need to be as consistent with QM as possible in

> > terms of

> > > numbers and units. Lets not get too concerned about cleaning up

> > every aspect of the risk parameters table to be a perfect match

> > with QM,

> > > though. Remember this is a guide of analyses to do and a

> > > compilation of

> > > screening values, not necessarily a formal spreadsheet work

> > > template> > (unless you guys have decided to do so??). Thanks to all,

> > > -Bob

>

> > \*\*\*\*\*

> > > Robert W. Gensemer, Ph.D.

> > > Parametrix, Inc.

> > > 33972 Texas Street SW

> > > Albany, OR 97321

> > > T 541-791-1667, x-6510

> > > F 541-791-1699

> > > C 541-760-1511

> > > rgensemer@parametrix.com

> > > \*\*\*\*\*

>

> > >>> <Benjamin.Shorr@noaa.gov> 1/10/2007 7:36:38 AM >>>

>

> > Eric-

>

> > > A few notes on the surface sediment screening numbers for

> > ecological> > risk:

>

> > > I strongly recommend that the units that are in this

> > > spreadsheet be

> > > changed to reflect the units in Query Manager. There should be a

> > > column

> > > with the units for each analyte (most metals in PPM, vols/svols

> > > etc> > PPB), and the guidelines should be adjusted to that for

> > > consistency.> >

> > > Total PCB's TEC should probably be .0598 (off by 10^3)

> > >

> > > Dieldrin (PPB) numbers are TEC/PEC = 1.9/61.8; spreadsheet has

> > > 2.85/6.7

> > >

> > > 2378 TCDD- there is one sample over 9 ng/kg (9E^-6 mg/kg) at 111

> > > under> railroad bridge. Looking directly at TCDD2378 conc. may

> > > benefit from

> > > a

> > > paired number.

> > >

> > > Hexachlorocyclohexane differs from QM TEC/PEC which is 2.37/4.99

> > > PPB,> spreadsheet has .94/1.38

> > >

> > > Hexachlorobutadiene, Tetrachloroethene, Trichloroethene units

> > > may be

> > > incorrect in spreadsheet (off by 10^3)

> > >

> > > Please let me know if there is a call today that I can join-  
> > otherwise> I'm available for the 1pm call tomorrow.  
> > Thanks,  
> > Ben  
> > >  
> > ----- Original Message -----  
> > From: Blischke.Eric@epamail.epa.gov  
> > Date: Tuesday, January 9, 2007 3:05 pm  
> > Subject: Re: A few modifications to the table  
> > >  
> > > Dana, here is a response to your questions and modifications to  
> > the> > table. I am copying the data evaluation folks and attaching  
> > your> > modifications to the table. I also have a few questions  
> > for Ben  
> > > regarding how QM handles certain summed values.  
> > >  
> > > I do not want to look at aluminum. 7600 mg/kg while screening  
> > in  
> > > at a  
> > > HQ of 0.1 is probably below background - upstream aluminum  
> > > concentrations range from 12,000 - 33,000 mg/kg. Further, the  
> > > direct  
> > > contact exposure scenarios are very conservative (350 days a  
> > year  
> > > for a  
> > > beach?).  
> > >  
> > > Regarding the TEQs and DDT, DDE and DDD sums - by manually, I  
> > > meant  
> > > thatit was not being calculated automatically by Query Manager.  
> > >  
> > > We  
> > > shouldbe able to do this in excel. I certainly support looking  
> > > at  
> > > the TEQs  
> > > but I want to get started on some easier evaluations first. We  
> > > may  
> > > haveto prioritize things here.  
> > >  
> > > Ben: What is included in the reported TEQ value - dioxin TEQs  
> > > or  
> > > dioxinand dioxin-like PCB TEQ?  
> > >  
> > > I don't really know how to best evaluate the PAHs. Regarding  
> > > naphthalene and Benzo(a)pyrene, we can look at these as  
> > individual> > chemicals. Hopefully,if we look at total PAHs,  
> > total low  
> > molecular> > weight PAHs and BAP and naphthalene, we will get a  
> > sense of the PAH  
> > > distribution to help us focus our evaluation. Another thing we  
> > might  
> > > want to do is query the carcinogenic PAHs and look at total  
> > > carcinogenicPAHs screened against BAP screening numbers.  
> > >  
> > > Ben: Do you know high molecular weight and low molecular weight  
> > > PAHs  
> > > are calculated.  
> > >  
> > > Regarding the modified table. I am ok with screening non-  
> > > carcinogens at  
> > > 0.1 (with the exception of Aluminum). Because QM is good at  
> > > looking at  
> > > concentration ranges, we should look at both HQ = 1 and HQ =  
> > 0.1.> >  
> > > I noticed the error regarding the residential soil PRG for BAP  
> > > (units  
> > > problem). You have correctly modified the screening number to  
> > be  
> > > 0.062mg/kg.  
> > >  
> > > Lets figure out the best way to too look at total PCBs (total  
> > > aroclors  
> > > or total congeners). For surface water, we should look at total  
> > > congeners due to interferences associated with the aroclor  
> > > results.  
> > > For  
> > > sediment, we should look at both total congeners and total  
> > aroclors.> > The total congeners represents a better number.  
> > However, we have  
> > much  
> > > less congener data than aroclor data. (PMX and Ben, I am  
> > > attaching a  
> > > write up on summing).  
> > >  
> > > Regarding TBT in Fish, our TBT data is limited to clams, and  
> > > juvenile  
> > > Chinook. Only one sample (a clam sample from the shipyard)  
> > > exceeds  
> > > thefish screening value (detected concentration = 530 ug/kg;  
> > fish  
> > > screeningnumber = 144 ug/kg; shellfish screening number = 1170  
> > > ug/kg). We can  
> > > still look at TBT in surface water.  
> > >  
> > > Eric  
> > >  
> > > (See attached file: 20070108Davoli Modif to ERIC  
> > > RiskParameters.xls)(Seeattached file: 20060201 Kissinger  
> > > Approach  
> > > Portland Harbor Upstream Fish  
> > > Tissue Sample Total PCBs, PCB TEQs, Dioxin\_Furan TEQs.doc)

[illegible]

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> > > for total low MW PAHs?
> > > 
> > > 
> > > I wasn't sure if TBT is above the SLV in fish. We can use the
> CRITFC> > Report value of 500 ug/mg as an SLV for all biota for
> lead but I
> don't
> know if we exceed this. For Hg in water, let's ues the ODEQ
> TMDL
> > > value.Ican look it up tomorrow.
> > > 
> > > I am in Health and Safety training on Tuesday but will try to
> call> > you
> > > at the morning break to discuss the table.
> > > 
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> > >
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